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ON A BASIS FOR "PEAKS OVER THRESHOLD" MODELING

by

M.R. Leadbetter



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Abstract. "Peaks over Threshold" ("POT") models commonly used e.g. in hydrology, assume that peak values of an iid or strionary sequence X<sub>i</sub> above a high value u, occur at Poisson points, and the exc. values of the peak above u are independent with an arbitrary common d.f. G. Motivation for these models has been provided by R.L. Smith (cf. [7],[8]), by using Pareto-type approximations of Pickands ([6]) for distributions of such excess values. These works strongly suggest that the Pareto family provides the appropriate class of distributions G for the POT model.

In the present paper we consider the point process of excess values of peaks above a high level u and demonstrate that this converges in distribution to a Compound Poisson Process as  $u \to \infty$  under appropriate assumptions. It is shown that the multiplicity distribution of this limit (i.e. the limiting distribution of excess values of peaks) must belong to the Pareto family and detailed forms are given for the normalizing constants involved. This exhibits the POT model specifically as a limit for the point process of excesses of peaks and delineates the distributions involved.

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## 1. Introduction.

In what are sometimes called "Peaks over Threshold" (POT) models (cf. [7]), the excess values over a high level u by an observed time series are assumed to occur at Poisson points and to have arbitrary common distributions. That is if  $\{X_i: i=1,2,\ldots\}$  is e.g. a stationary (or iid) sequence, the exceedance points  $\{i: X_i > u\}$  are assumed to be Poisson and the corresponding excess values  $(X_i-u)_+$  to be independent with an arbitrary distribution.

The Poisson nature of the occurrence of exceedances is intuitively clear since e.g. if  $X_i$  are iid with d.f. F, the number of exceedances of a high level  $u_n$  by  $X_1, \ldots, X_n$ , is binomial with parameters  $(n, 1-F(u_n))$  and hence approximately Poisson if  $n(1-F(u_n))$  converges to some value  $\tau > 0$ . This will be made more precise below by a time normalization. Further motivation for the model is provided by R.L. Smith ([7],[8]) based on theory of Pickands [6], restricting the distribution for excess values to a "generalized Pareto" (GP) form

(1.1) 
$$G_{\alpha,\beta}(x) = 1 - (1 + \alpha x/\beta)^{-1/\alpha} \quad \beta > 0, \ \alpha \neq 0$$
$$= 1 - e^{-x/\beta} \qquad \beta > 0, \ \alpha = 0$$

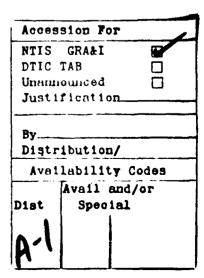
where the range of x is  $(0,\infty)$  if  $\alpha \ge 0$  and  $(0,-\alpha^{-1}\beta)$  if  $\alpha < 0$ .

This class is a flexible 2-parameter family, but more importantly for a wide class of F's the excess distribution

(1.2) 
$$F_u(x) = P\{X - u \le x | X > u\}$$

is approximately GP in a sense shown by Pickands [6], viz.

(1.3) 
$$\inf_{\beta} \sup_{\mathbf{x}} |F_{\mathbf{u}}(\mathbf{x}) - G_{\alpha,\beta}(\mathbf{x})| \to 0 \quad \text{as } \mathbf{u} \to \infty$$



for some fixed  $\alpha$ . That is for high levels u,  $\beta$  may be chosen (depending on u) so that  $G_{\alpha,\beta}(x)$  is uniformly close to  $F_{u}(x)$ .

Write  $x_F$  for the right endpoint (sup{x: F(x) < 1} of the d.f. F and  $\overline{F}(x) = 1 - F(x)$ , the tail of F. Then the class of d.f.'s F for which the above GP approximation holds includes all those satisfying

(1.4) 
$$\overline{F}(u+xg(u))/\overline{F}(u) \rightarrow \overline{G}(x)$$
 as  $u \rightarrow x_F$ 

for some function g(u) > 0, some d.f. G and all  $0 < x < x_F (\le \infty)$ . It is known ([6]) that any such G in (1.4) must be G.P. as in (1.1) for some  $\alpha, \beta$ .

Note that (1.4) holds for all d.f.'s F of interest in extreme value theory i.e. such that if  $M_n = \max(X_1, \dots, X_n)$ ,  $P\{a_n(M_n - b_n) \le x\}$  (=  $F^n(x/a_n + b_n)$ ) has a non degenerate limit  $\Lambda(x)$ . For example if  $\overline{G}(x) = e^{-x}$ , (1.1) is a classical domain of attraction criterion for a "Type I" extreme value distribution  $\Lambda(x) = \exp(-e^{-x})$ . If instead F has a regulary varying tail  $(1-F(ux))/(1-F(u)) \to x^{-\alpha}$ ,  $\alpha>0$ , as  $u \to \infty$ , each x>0, then (1.1) holds with  $\overline{G}(x) = (1+x)^{-\alpha}$ . g(u) = u, and  $\Lambda(x)$  is then "Type II" i.e.  $\Lambda(x) = \exp(-x^{-\alpha})$ , x>0.

Hence in a wide variety of cases of interest the distribution  $F_u(x)$  of excesses (given by (1.2)) of the level u is approximately GP,  $G_{\alpha,\beta}(x)$  in the sense stated, where  $\alpha$  is fixed but  $\beta$  can change with u. As discussed in [7] this provides significant intuitive support for the POT model. Our purpose here is to further justify the model by exhibiting it as the limit of point processes of excesses of high levels, the limiting distribution of excess values being shown to be GP,  $G_{\alpha,\beta}(x)$  where now  $\beta$  as well as  $\alpha$ , is independent of u.

For clarity this will be shown for iid sequences in Section 2 and extended to dependent (mixing) situations in Section 3. In the latter case high serial correlation can cause clustering of exceedances of high levels and the peak values are then defined to be the largest in each cluster. Two points worth noting are (i) the dependence modifies the theory via the introduction of a single parameter, the "extremal index" (essentially the inverse of mean cluster size) and (ii) the GP form applies to the peak values but not necessarily to other cluster properties such as their lengths. A corresponding theory will be indicated in Section 3 for other such cases.

## 2. The iid case.

Let  $\{u_n\}$  be a sequence of levels such that

(2.1) 
$$n(1-F(u_n)) \rightarrow \tau > 0.$$

Define point processes  $N_n$  to consist of the points i/n for which  $X_i > u_n$ , i.e. the exceedance points normalized by 1/n.  $N_n$  is thus defined on the positive real line but it will be convenient to restrict attention to the unit interval, corresponding to exceedances among the n sample values  $X_1, \ldots, X_n$ . Hence  $N_n(B)$  is defined for (Borel) subsets  $B \subset (0,1]$  by  $N_n(B) = \#\{i/n \in B: X_i > u_n\}$ .

It is trivial to show that under (2.1) that  $N_n \xrightarrow{d} N$  where N is a Poisson Process on (0,1] with intensity  $\tau$ . For if I = (a,b]  $\subset$  (0,1],  $N_n(I)$  is binomial with parameters [nb]-[na],  $p_n=1-F(u_n)$  ([] denoting integer part) and  $np_n \to \tau$  so that

$$P\{N_n(I) = r\} \longrightarrow e^{-\tau(b-a)} [\tau(b-a)]^{r}/r! = P\{N(I) = r\}.$$

Hence  $N_n(I) \xrightarrow{d} N(I)$  and by independence if  $I_1, \dots, I_k$  are disjoint

$$(N_n(I_1), \dots N_n(I_k)) \xrightarrow{d} (N(I_1), \dots N(I_k))$$

which is sufficient ([2]) to show full weak convergence  $N_n \xrightarrow{d} N$ .

Now associate with each point of N  $_n$  (i/n such that X  $_i$  > u  $_n$ ) the corresponding excess value X  $_i$ -u  $_n$  to give the "point process N  $_n^*$  of excesses".

Technically perhaps  $N_n^*$  should be regarded as a "marked point process" (or as an atomic random measure) since the  $(X_i - u_n)$  are not necessarily integer valued but it will be convenient (and legitimate) to call it a point process whose events (at  $(i/n: X_i > u_n)$ ) have (not necessarily integer) multiplicities  $(X_i - u_n)$ .

Since, as above, the positions of the excesses converge to a Poisson Process, it seems evident that one should expect any limit  $N^*$  for  $N^*$  to be a compound Poisson process having events at Poisson points with intensity  $\tau$  and (independent) multiplicities with some common d.f. G. Such a result may be shown (cf. [1]) but here we give sufficient conditions for such convergence. Here and below we write  $N^* = CP(\tau,G)$  to denote a Compound Poisson point process whose Poisson events have intensity  $\tau$  with (not necessarily integer-valued) multiplicities having d.f. G.

Theorem 2.1. Let  $X_i$ , i=1,2... be iid with d.f. F satisfying (1.4) for some g. G, and let  $(u_n)$  be levels satisfying (2.1). Then  $a_n \stackrel{\times}{n} \xrightarrow{d} \stackrel{\times}{N}$ ,  $CP(\tau,G)$  where  $\tau$  is as in (2.1), G as in (1.4) and  $a_n = 1/g(u_n)$ .

<u>Proof</u>: If  $G_n(x)$  (=  $F_{u_n}(x/a_n)$ ) denotes the conditional d.f. of  $a_n(X_1-u_n)$  given  $X_1 > u_n$ , then clearly

$$exp\{-sa_n(X_1-u_n)_+\} = F(u_n) + \overline{F}(u_n) \int_0^\infty e^{-sx} dG_n(x).$$

But from (1.4),  $G_n(x) \to G(x)$  and hence  $\int_0^\infty e^{-sx} dG_n(x) \to \phi(s) = \int_0^\infty e^{-sx} dG(x)$  so that if  $I = (a,b] \subset (0,1]$  contains  $m_n$  points i/n  $(m_n \sim n(b-a))$ 

$$\mathcal{E}^{m} \exp\{-sa_{n}(X_{1}-u_{n})_{+}\} = [1-(\tau/n)(1-\phi(s))(1+o(1))]^{m}n$$

$$\to e^{-\tau(b-a)(1-\phi(s))}$$

which is the Laplace Transform  $e^{-sN^*(a,b]}$  where  $N^*$  is  $CP(\tau,G)$ . Since  $N_n^*(I)$  is

the sum  $\sum_{i/n\in I} (X_i - u_n)_+$  of  $m_n$  iid terms, it follows that  $a_n N_n^*(I) \xrightarrow{d} N^*(I)$ . Hence by independence

$$(\mathbf{a_n} \mathbf{N_n^{\star}}(\mathbf{I_1}) \dots \mathbf{a_n} \mathbf{N_n^{\star}}(\mathbf{I_n})) \xrightarrow{\mathbf{d}} (\mathbf{N_n}(\mathbf{I_1}) \dots \mathbf{N_n}(\mathbf{I_k}))$$

for any disjoint  $I_1, \ldots, I_k$  so that  $a_n \stackrel{*}{N_n} \stackrel{d}{\longrightarrow} \stackrel{N}{N}$  as required.

Thus in the iid case the (normalized) point process  $a_n N_n^{\bowtie}$  of excesses over  $u_n$  has a Compound Poisson limit and may be regarded as approximately  $CP(\tau,G)$  for large n. Note again that G is a GP distribution  $G_{\alpha,\beta}$  for some fixed  $\alpha,\beta$ . This provides a strong basis for the POT model in the iid case.

# 3. Stationary sequences

A stationary sequence  $X_1, X_2, \ldots$  can exhibit both "long range" and "local" dependence between the  $X_i$ . Here the former, but not the latter, will be restricted by a "mixing condition" which enables the collection of  $X_i$ 's into groups which are approximately independent but with possible high dependence within groups. Strong mixing will suffice to restrict long range dependence. However one may "tailor" this to the problem at hand with a slightly weaker restriction  $\Delta(v_n)$  defined for any sequence of constants  $v_n$  as follows. For  $1 \le j \le k \le n$  write  $\mathfrak{B}_{jk}(v_n) = \sigma\{(X_s - v_n)_+ : j \le s \le k\}$  and say that  $\Delta(v_n)$  holds if for  $1 \le \ell \le n$  whenever  $A \in \mathfrak{B}_{1,j}(v_n)$ ,  $B \in \mathfrak{B}_{j+\ell,n}(v_n)$ ,  $1 \le j \le n-\ell$  and  $\alpha_{n,\ell} \to 0$  for some  $\ell_n = o(n)$ .

High <u>local</u> dependence is reflected in the presence of clustering of exceedances of high levels. "Clusters" will be defined precisely below, but we first note that for levels  $(u_n)$  satisfying (2.1) the mean size of a cluster typically converges to a parameter whose inverse  $\theta$  is sometimes called the "extremal index" of the sequence  $\{X_i\}$ . Specifically  $\theta$  is defined by the property that if  $M_n = \max(X_1, X_2, \dots, X_n)$  and  $u_n$  satisfies (2.1) then

$$(3.1) P(M_n \le u_n) \to e^{-\theta \tau}.$$

 $\theta$  being independent of  $\tau$ . (For i.i.d. sequences  $\theta = 1$ ). Such  $\theta$  exists under general conditions (cf [4], Sec. 3.7)

Clusters of exceedances of high levels may be defined in various ways, the most obvious being as runs of consecutive exceedances. However if  $\Delta(v_n)$  holds for a given sequence  $\{v_n\}$  of levels the following "block definition" is more convenient (and often asymptotically equivalent to that using runs). Choose integers  $k_n \to \infty$ ,  $k_n = o(n)$ , satisfying

(3.2) 
$$k_n(\alpha_n, \ell_n + \ell_n/n) \to 0.$$

Write  $r_n = [n/k_n]$  and divide the integers (1...n) into consecutive blocks

$$J_{i} = \{(i-1)r_{n}+1, (i-1)r_{n}+2, \dots, ir_{n}\} \quad 1 \le i \le k_{n}$$

$$J_{k_{n}+1} = \{k_{n}r_{n}+1, k_{n}r_{n}+2, \dots, n\}$$

and regard the exceedances (if any) in a block as forming a cluster. The choice of  $k_n$  (or equivalently  $r_n$ ) is flexible, subject to the growth restriction (3.2).

Obviously this block definition can count a run of consecutive exceedances as two or more clusters, if the run straddles more than one block so that the block definition is less natural in some cases. However it can be more appropriate than the runs definition for sequences with high local variability. In any case we use the block definition for its convenience. Further if the block  $J_i$  contains exceedances the first point,  $(i-1)r_n+1$ , of  $J_i$  will be regarded as the <u>location</u> of the cluster in  $J_i$ .

Define now a point process  $P_n$  of normalized cluster positions, i.e. consisting of the points  $\{((i-1)r_n+1)/n, 1 \le i \le k_n : M(J_i) > u_n\}$  where M(E) is

written for  $\max\{X_j\colon j\epsilon E\}$ . Associate with each point  $((i-1)r_n+1)/n$  of  $P_n$  the corresponding maximum excess  $M(J_i)-u_n=\max\{(X_j-u_n)_+\colon j\in J_i\}$ , giving the (again technically "marked") point process  $P_n^*$  of peak excess values above the level in the clusters. We show that a result like Theorem 2.1 holds in the dependent case with  $P_n^*$  replacing  $N_n^*$ . In fact it may be shown (cf [5]) that for i.i.d. sequences  $N_n$  and  $P_n$  are asymptotically equivalent in a strong sense as also are  $N_n^*$  and  $P_n^*$  so that  $P_n$  and  $P_n^*$  generalize  $N_n$  and  $N_n^*$ .

It was shown in Section 1 that in the iid case the exceedance point process  $N_n$  has a Poisson limit with intensity  $\tau$ . This result is simply generalized (cf. [1]) to show that under dependence  $P_n$  has a Poisson limit (with intensity  $\theta\tau$ ) whereas  $N_n$  itself has a Compound Poisson limit whose events occur a\* the positions of clusters (i.e. points of  $P_n$ ), with multiplicities given by (limiting) cluster sizes. However the (limiting) distribution for multiplicities need not have a GP form and need not be totally determined by the (tail of the) marginal d.f. F of the  $X_i$ . On the other hand it will be shown below that  $P_n^*$  has a CP( $\theta\tau$ ,G) distributional limit where G is obtained from the tail of F via (1.4). Thus G has GP form and the dependence influences the limit only through the factor  $\theta$  in the intensity of the underlying Poisson Process.

The Compound Poisson limit for  $P_n^*$  will be obtained by considering the asymptotic behavior of the maximum in a cluster and showing that the clusters are essentially independent. The specific basic results needed are contained in the following lemma.

Lemma 3.1. Let  $\{X_n\}$  be stationary with extremal index  $\theta > 0$  and marginal d.f. F satisfying (1.4), and let  $\Delta(v_n)$  hold with  $v_n = u_n + xg(u_n)$ , each  $x \ge 0$ , where  $u_n$  satisfies (2.1). Let  $\{k_n\}$  satisfy (3.2) and write  $a_n = (g(u_n))^{-1}$ . Then (with  $v_n = [n/k_n]$ ), as  $n \to \infty$ 

(i) 
$$P\{a_n(M_{r_n}-u_n) > x\} \sim \frac{\theta \tau}{k_n} \overline{G}(x), x \ge 0$$

(ii) 
$$P\{a_n(M_{r_n}-u_n) \le x | M_{r_n}>u_n\} \longrightarrow G(x)$$

$$\text{(iii) } \left[ \exp \left\{ -\operatorname{sa}_{n} \left( \operatorname{M}_{r_{n}} - \operatorname{u}_{n} \right)_{+} \right\} \right]^{k_{n}} \longrightarrow \exp \left\{ -\theta \tau \left( 1 - \phi(s) \right) \right\}$$

where 
$$\phi(s) = \int_0^\infty e^{-sx} dG(x)$$

Proof: Since  $n[1-F(u_n^{+a_n^{-1}x})] \sim \tau \overline{F}(u_n^{+xg}(u_n^{-1}))/\overline{F}(u_n^{-1}) \to \tau \overline{G}(x)$  and  $\{X_n^{-1}\}$  has extremal index  $\theta$ , it follows that

$$P\{M_n \le u_n + a_n^{-1}x\} \longrightarrow \exp(-\theta \tau \overline{G}(x)).$$

But it follows in a standard way from the mixing condition  $\Delta(v_n)$  (cf. [1, Lemma 2.3]) that  $P\{M_n \le v_n\} - P^{k_n}\{M_r \le v_n\} \to 0$  so that  $P^{k_n}\{M_r \le u_n + a_n^{-1}x\} \to \exp(-\theta \tau \overline{G}(x))$  from which (i) readily follows. The left hand side of (ii) is

1 - 
$$P\{M_{r_n} > u_n + a_n^{-1}x\} / P\{M_{r_n} > u_n\} \approx 1 - \frac{\theta \tau}{k_n} \overline{G}(x) / \frac{\theta \tau}{k_n} (1 + o(1))$$

by (i), so that (ii) follows. Finally (iii) follows by the first calculation of Theorem 2.1 with M<sub>r</sub> replacing  $X_i$ , using (ii).

The main result now follows in a similar way to Theorem 2.1 on using approximate independence between the clusters.

Theorem 3.2. Let  $\{X_n\}$  be stationary with extremal index  $\theta > 0$ , and marginal d.f F satisfying (1.4), and let  $\Delta(v_n)$  hold with  $v_n = u_n + xg(u_n)$  each  $x \ge 0$ , where  $u_n$  satisfies (2.1). Let  $k_n \to \infty$  satisfy (3.2),  $r_n = \lfloor n/k_n \rfloor$  and  $a_n = (g(u_n))^{-1}$ . Then the point process  $P_n^*$  of peak values above  $u_n$  satisfies  $a_n P_n^* \xrightarrow{d} P_n^*$  where  $P_n^*$  is  $CP(\theta \tau, G)$  and G is as in (1.4).

Proof: Let I be a subinterval of (0,1] and  $J_i = \{(i-1)r_n+1,\ldots,ir_n\}$  as defined above,  $1 \le i \le k_n$ . Then it follows from the mixing conditions along the same lines as Lemma 2.2 of [3] (or Lemma 2.2 of [1]) that

$$\operatorname{\operatorname{\mathcal{E}exp}} \{-\operatorname{sa}_n\operatorname{P}_n^{\bigstar}(\operatorname{I})\} - \operatorname{\operatorname{II}} \operatorname{\operatorname{\mathcal{E}exp}} \{-\operatorname{sa}_n\operatorname{P}_n^{\bigstar}(\operatorname{J}_i)\} \longrightarrow 0$$

as  $n \to \infty$ , so that

$$\begin{aligned} \exp\{-\operatorname{sa}_{n}\operatorname{P}_{n}^{*}(I)\} &= \left[\operatorname{\operatorname{\mathcal{E}exp}}\{-\operatorname{sa}_{n}(\operatorname{M}_{\Gamma_{n}}^{-\operatorname{u}}_{n})_{+}\}\right]^{k_{n}}\operatorname{m}(I)(1+o(1)) \\ & \longrightarrow \exp\{-\theta\tau\operatorname{m}(I)(1-\phi(s))\} \end{aligned}$$

by Lemma 3.1, where  $\phi(s) = \int_0^\infty e^{-sx} dG(x)$ . Hence  $a_n P_n^*(I) \xrightarrow{d} P^*(I)$  where  $P^*$  is  $CP(\theta\tau,G)$  on (0,1]. Now if  $I_1,\ldots,I_k$  are disjoint subintervals of (0,1] it follows also as in Lemma 2.2 of [3] (or Lemma 2.2 of [1]) that

$$\operatorname{\operatorname{\mathcal{E}exp}} \{-a_n \sum_{j=1}^k s_j P_n^{\bigstar}(I_j)\} - \prod_{j=1}^k \operatorname{\operatorname{\mathcal{E}exp}} \{-a_n P_n^{\bigstar}(I_j)\} \longrightarrow 0$$

from which it follows that

$$(a_n P_n^*(I_1) \dots a_n P_n^*(I_k)) \xrightarrow{d} (P^*(I_1) \dots P^*(I_k))$$

and hence that  $a_n \stackrel{*}{P} \xrightarrow{d} \stackrel{*}{P}$ .

Finally we reiterate that this result is one of many which can be obtained involving different aspects of cluster structure. For example the Compound Poisson limit for  $N_n$  was cited above, the multiplicities corresponding to cluster sizes. More complicated functions — such as the sum of powers of excess values in a cluster — may also be considered and will lead to Compound Poisson limits. Such cases may be useful in applications where damage from high levels (e.g. high pollution episodes) may be modeled as a specific

function of the excess values. However the case of (excess) <u>peak</u> values in a cluster is especially important (e.g. in describing severe floods, where damage can well be a function of flood level). Theorem 3.2 establishing the POT approximation is particularly useful since the multiplicity distribution then depends only on the marginal d.f. F and moreover is known to have GP form.

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